=> fil req

FILE 'REGISTRY' ENTERED AT 07:37:39 ON 01 APR 2009

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STRUCTURE FILE UPDATES: 30 MAR 2009 HIGHEST RN 1129871-47-1 DICTIONARY FILE UPDATES: 30 MAR 2009 HIGHEST RN 1129871-47-1

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH January 9, 2009.

Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

http://www.cas.org/support/stngen/stndoc/properties.html

=> d 120 ide can tot

- L20 ANSWER 1 OF 5 REGISTRY COPYRIGHT 2009 ACS on STN
- RN 773889-61-5 REGISTRY
- ED Entered STN: 02 Nov 2004
- CN Benzamide, N, N'-1, 4-phenylenebis[3,5-diamino- (CA INDEX NAME)
- MF C20 H20 N6 O2
- SR CA
- LC STN Files: CA, CAPLUS, CASREACT

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

2 REFERENCES IN FILE CA (1907 TO DATE)
2 REFERENCES IN FILE CAPLUS (1907 TO DATE)

REFERENCE 1: 146:462594

REFERENCE 2: 141:332962

- L20 ANSWER 2 OF 5 REGISTRY COPYRIGHT 2009 ACS on STN
- RN 773889-59-1 REGISTRY
- ED Entered STN: 02 Nov 2004
- CN Benzamide, N,N'-(oxydi-4,1-phenylene)bis[3,5-diamino- (9CI) (CA INDEX

NAME)

MF C26 H24 N6 O3

CI COM

SR CA

LC STN Files: CA, CAPLUS

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

REFERENCE 1: 141:332962

- L20 ANSWER 3 OF 5 REGISTRY COPYRIGHT 2009 ACS on STN
- RN 773889-55-7 REGISTRY
- ED Entered STN: 02 Nov 2004
- CN Methanone, 1,1'-(1,4-piperazinediyl)bis[1-(3,5-diaminophenyl)- (CA INDEX NAME)

OTHER CA INDEX NAMES:

- CN Piperazine, 1,4-bis(3,5-diaminobenzoyl)- (9CI)
- MF C18 H22 N6 O2
- CI COM
- SR CA
- LC STN Files: CA, CAPLUS

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1907 TO DATE) 1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

REFERENCE 1: 141:332962

- L20 ANSWER 4 OF 5 REGISTRY COPYRIGHT 2009 ACS on STN
- RN 518992-19-3 REGISTRY
- ED Entered STN: 22 May 2003
- CN 1,3-Benzenediamine, 5,5'-(sulfonyldi-5,2-benzoxazolediyl)bis- (9CI) (CA INDEX NAME)
- MF C26 H20 N6 O4 S
- CI COM
- SR Chemical Library

Supplier: Ambinter LC STN Files: CA, CAPLUS, CHEMCATS

$$\begin{array}{c} \text{NH2} \\ \text{H2N} \end{array}$$

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1907 TO DATE) 1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

REFERENCE 1: 141:332962

L20 ANSWER 5 OF 5 REGISTRY COPYRIGHT 2009 ACS on STN

176258-99-4 REGISTRY RN

ED Entered STN: 15 May 1996

9H-Fluorene-2,7-diamine, 9,9-bis(4-aminophenyl)- (CA INDEX NAME) CN

OTHER NAMES: CN 2,7-Diamino-9,9-bis(4-aminophenyl)fluorene

MF C25 H22 N4 COM

CI SR CA

LC STN Files: CA, CAPLUS, CASREACT

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

8 REFERENCES IN FILE CA (1907 TO DATE)

1 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

8 REFERENCES IN FILE CAPLUS (1907 TO DATE)

REFERENCE 1: 141:332962

REFERENCE 2: 139:181414

REFERENCE 3: 134:179710

REFERENCE 134:179368 4:

REFERENCE 5: 131:102606

REFERENCE 6: 130:125670

REFERENCE 7: 126:277826

REFERENCE 8: 124:316706

=> d sta que 124 L22 S

NODE ATTRIBUTES: DEFAULT MLEVEL IS ATOM DEFAULT ECLEVEL IS UNLIMITED

GRAPH ATTRIBUTES: RSPEC 15 25 17 32 NUMBER OF NODES IS 30

100.0% PROCESSED 98 ITERATIONS SEARCH TIME: 00.00.01 1 ANSWERS

=> d ide can 124

L24 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2009 ACS on STN

RN 168914-99-6 REGISTRY

ED Entered STN: 13 Oct 1995

CN 5,6'-Bibenzoxazole, 2,2'-bis(4-fluorophenyl)- (CA INDEX NAME)

MF C26 H14 F2 N2 O2

SR CA

LC STN Files: CA, CAPLUS



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

REFERENCE 1: 123:229141

=> fil hcaplus

FILE 'HCAPLUS' ENTERED AT 07:37:57 ON 01 APR 2009
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FILE COVERS 1907 - 1 Apr 2009 VOL 150 ISS 14 FILE LAST UPDATED: 31 Mar 2009 (20090331/ED)

HCAplus now includes complete International Patent Classification (IPC) reclassification data for the third quarter of 2008.

CAS Information Use Policies apply and are available at:

http://www.cas.org/legal/infopolicy.html

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> d 134 bib abs hitstr tot

L34 ANSWER 1 OF 14 HCAPLUS COPYRIGHT 2009 ACS on STN AN 2008:1534289 HCAPLUS Full-text

DN 150:57016

TI Manufacture of 6,6-polyimide copolymer soluble in solvent IN Itatani, Biroshi

IN Itatani, Hiroshi

PA Solpit Industries, Ltd., Japan; Sojitz Corporation

SO PCT Int. Appl., 35pp. CODEN: PIXXD2

DT Patent LA Japanese

FAN.CNT 1

FAN.	CNT	1																	
	PATENT NO.					KIND		DATE		APPLICATION NO.						DATE			
PI	WO 2008155811				A1		20081224		WO 2007-JP62247					20070618					
		w:	ΑE,	AG,	AL,	AM,	AT,	AU,	AZ,	BA,	BB,	BG,	BH,	BR,	BW,	BY,	BZ,	CA,	
			CH,	CN,	CO,	CR,	CU,	CZ,	DE,	DK,	DM,	DO,	DZ,	EC,	EE,	EG,	ES,	FI,	
			GB,	GD,	GE,	GH,	GM,	GT,	HN,	HR,	HU,	ID,	IL,	IN,	IS,	JP,	KE,	KG,	
			KM,	KN,	KP,	KR,	ΚZ,	LA,	LC,	LK,	LR,	LS,	LT,	LU,	LY,	MA,	MD,	MG,	
			MK,	MN,	MW,	MX,	MY,	ΜZ,	NA,	NG,	NI,	NO,	ΝZ,	OM,	PG,	PH,	PL,	PT,	
			RO,	RS,	RU,	SC,	SD,	SE,	SG,	SK,	SL,	SM,	SV,	SY,	TJ,	TM,	TN,	TR,	
			TT,	TZ,	UA,	UG,	US,	UZ,	VC,	VN,	ZA,	ZM,	ZW						
		RW:	ΑT,	BE,	BG,	CH,	CY,	CZ,	DE,	DK,	EE,	ES,	FΙ,	FR,	GB,	GR,	HU,	IE,	
			IS,	IT,	LT,	LU,	LV,	MC,	MT,	NL,	PL,	PT,	RO,	SE,	SI,	SK,	TR,	BF,	
			ВJ,	CF,	CG,	CI,	CM,	GA,	GN,	GQ,	GW,	ML,	MR,	ΝE,	SN,	TD,	TG,	BW,	
			GH,	GM,	KE,	LS,	MW,	MZ,	NA,	SD,	SL,	SZ,	TZ,	UG,	ZM,	ZW,	AM,	AZ,	

BY, KG, KZ, MD, RU, TJ, TM

PRAI WO 2007-JP62247 20070618

B The polyimide is synthesized from 6,6-imide segment which is an imide oligomer having pyromellitic dianhydride (PMDA) at both ends produced by adding 4 molar equiv of PMDA and 2 molar equiv of diaminotoluene (DAT) to an imide oligomer produced by heating 1 molar equiv of biphenyltetracarboxylic dianhydride (BPDA) and 2 molar equiv of diaminodiphenyl ether (DADE) to 160 to 200° in an organic polar solvent in the presence of a catalyst.

T 1093221-10-3P 1093221-11-4P 1093221-12-5P

1093221-13-6P

RL: IMF (Industrial manufacture); PRP (Properties); PREP (Preparation) (manufacture of 6,6-polyimide copolymer soluble in solvent)

RN 1093221-10-3 HCAPLUS CN 1H.3H-Benzoll.2-c:4.5

1H,3H-Benzo[1,2-c:4,5-c']difuran-1,3,5,7-tetrone, polymer with 3-(4-aminophenoxy)benzenamine, [5,5'-biisobenzofuran]-1,1',3,3'-tetrone, 1,4-bis(3,5-diaminobenzoy1)piperazine and ar-methyl-1,3-benzenediamine, block (CA INDEX NAME)

CM 1

CRN 773889-55-7 CMF C18 H22 N6 O2

CM 2

CRN 26764-44-3 CMF C7 H10 N2 CCI IDS

D1-Me

CM 3

CRN 2657-87-6 CMF C12 H12 N2 O

CM 4

CRN 2420-87-3 CMF C16 H6 O6

CM 5

CRN 89-32-7 CMF C10 H2 O6

RN 1093221-11-4 HCAPLUS

1H,3H-Benzo[1,2-c:4,5-c']difuran-1,3,5,7-tetrone, polymer with 3-(4-aminophenoxy) benzenamine, [5,5'-biisobenzofuran]-1,1',3,3'-tetrone, ar-methyl-1,3-benzenediamine and 1,1'-(1,4-piperazinediy1) bis[1-(3,5-diaminophenyl)methanone], block (CA INDEX NAME)

CM 1

CN

CRN 773889-55-7 CMF C18 H22 N6 O2

CM 2

CRN 97917-34-5 CMF (C2 H6 O Si)n C10 H28 N2 O Si2 CCI PMS

CM 3

CRN 26764-44-3 CMF C7 H10 N2 CCI IDS

D1—Me

CM 4

CRN 2657-87-6 CMF C12 H12 N2 O

CM 5

CRN 2420-87-3 CMF C16 H6 O6

CM 6

CRN 89-32-7 CMF C10 H2 O6

RN 1093221-12-5 HCAPLUS

CN 1H,3H-Benzo[1,2-c:4,5-c']difuran-1,3,5,7-tetrone, polymer with $\alpha = [(3-\min n propyl) dimethylsily1]-\omega = [((3-\min n propyl) dimethylsily1) ewy [((3-\min n propyl) dimethylsily1) ewy [(3-min propyl) dimethylsily1) end [(5,5'-biisobenzofuran]-1,1',3,3'-tetrone, 1,4-bis(3,5-diaminobenzoyl)piperazine and ar-methyl-1,3-benzenediamine, block (CA INDEX NAME)$

CM 1

CRN 773889-55-7 CMF C18 H22 N6 O2

CM 2

CRN 97917-34-5

CMF (C2 H6 O Si)n C10 H28 N2 O Si2

CCI PMS

CM 3

CRN 26764-44-3

CMF C7 H10 N2

CCI IDS

D1-Me

CM 4

CRN 2420-87-3

CMF C16 H6 O6

CM 5

CRN 89-32-7 CMF C10 H2 O6

RN 1093221-13-6 HCAPLUS

CN 1H,3H-Benzo[1,2-c:4,5-c']difuran-1,3,5,7-tetrone, polymer with [5,5'-biisobenzofuran]-1,1',3,3'-tetrone, 1,4-bis(3,5-diaminobenzoy]lpiperazine, ar-methyl-1,3-benzenediamine and

3,3'-[1,4-phenylenebis(oxy)]bis[benzenamine], block (CA INDEX NAME)

CM :

CRN 773889-55-7 CMF C18 H22 N6 O2

11

CM 2

CRN 59326-56-6 CMF C18 H16 N2 O2

CM 3

CRN 26764-44-3 CMF C7 H10 N2 CCI IDS

D1-Me

CM 4

CRN 2420-87-3 CMF C16 H6 O6

CM 5

CRN 89-32-7 CMF C10 H2 O6

RE.CNT 17 THERE ARE 17 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT

L34 ANSWER 2 OF 14 HCAPLUS COPYRIGHT 2009 ACS on STN

AN 2007:240212 HCAPLUS Full-text

DN 146:462594

TI Facile Synthesis of Amine-Terminated Aromatic Polyamide Dendrimers via a Divergent Method

AU Washio, Isao; Shibasaki, Yuji; Ueda, Mitsuru

CS Department of Organic and Polymeric Materials, Graduate School of Science and Engineering, Tokyo Institute of Technology, 2-12-1-H120 O-okayama, Meguro-ku, Tokyo, 152-8552, Japan

SO Organic Letters (2007), 9(7), 1363-1366

CODEN: ORLEF7; ISSN: 1523-7060 PB American Chemical Society

DT Journal

LA English

OS CASREACT 146:462594

AB A novel, rapid, inexpensive, and highly efficient divergent approach for the synthesis of a 32-amine-terminated 64 polyamide dendrimer has been developed. Each generation dendrimer was successfully obtained by the condensation of the preceding generation dendrimer with the building block and the deprotection with hydrazine in one pot. All the dendrimers were easily purified by precipitation in alkaline water, and the purity was confirmed by NMR, MALDITOF mass spectra, and elemental anal.

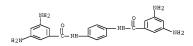
IT 773889-61-5P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(facile synthesis of amine-terminated aromatic polyamide dendrimers via divergent method)

RN 773889-61-5 HCAPLUS

CN Benzamide, N, N'-1, 4-phenylenebis[3, 5-diamino- (CA INDEX NAME)



RE.CNT 36 THERE ARE 36 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT

L34 ANSWER 3 OF 14 HCAPLUS COPYRIGHT 2009 ACS on STN

AN 2004:847588 HCAPLUS Full-text

DN 141:332962

TI Crosslinked polyimides, compositions containing them and method for their manufacture

IN Itatani, Hiroshi

PA Fi R & D Co. Ltd., Japan

SO PCT Int. Appl., 68 pp.

CODEN: PIXXD2

DT Patent

LA Japanese

FAN.CNT 1

PATENT NO. KIND DATE APPLICATION NO. DATE

```
PΤ
    WO 2004087793
                         A1
                               20041014 WO 2004-JP4305
                                                                   20040326 <--
         W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH,
             CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD,
             GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC,
             LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI,
             NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY,
             TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW
         RW: BW, GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ,
            BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE,
             ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PL, PT, RO, SE, SI,
             SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN,
             TD, TG
     EP 1614704
                               20060111
                                         EP 2004-723799
                                                                   20040326 <--
                         A1
        R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,
             IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, PL, SK
     US 20070106056
                         A1
                               20070510
                                           US 2005-550887
                                                                   20050228 <--
PRAI JP 2003-90546
                         А
                               20030328 <--
                               20030417 <--
     JP 2003-112425
                         Α
     JP 2003-412832
                         Α
                               20031211 <--
     WO 2004-JP4305
                         W
                               20040326
                                         <--
     The crosslinked polyamides are produced by the polycondensation of a
AB
     tetraamine, a tetracarboxylic acid dianhydride and an aromatic diamine in the
     presence of a catalyst. The crosslinked polyamides exhibit a dielec. constant
     of ≤2.7 while compns. containing polyimides have inherent good heat
     resistance, elec. insulation and chemical resistance, and are useful for elec.
     and electronic device manufacture Thus, polycondensing bis(3,5-
     diaminobenzoyl)-1,4-piperazine with biphenyltetracarboxylic dianhydride and
     4.4'-diaminodiphenvl ether using oxalic acid and pyridine 2 component catalyst
     in N-methyl-2-pyrrolidone then coupling with 3,3',4,4'-diphenyl ether
     tetracarboxylic dianhydride and 1,3-bis(4-aminophenyl)benzene gave a
     crosslinked polyimide having the claimed properties.
     773889-56-8p 773889-57-9p 773889-58-0p
     773889-62-6P 773889-63-7P 773889-64-8P
     773889-65-9P 773889-66-0P 773889-67-1P
     773889-68-2P 773889-69-3P 773889-70-6P
     773889-71-7p 773889-72-8p 773889-73-9p
     773889-75-1P 773889-76-2P 773889-77-3P
     RL: IMF (Industrial manufacture); PRP (Properties); TEM (Technical or
     engineered material use); PREP (Preparation); USES (Uses)
        (crosslinked polyimides with low dielec. constant, compns. containing them
        and method for their manufacture and use)
    773889-56-8 HCAPLUS
RN
CN
    [5,5'-Biisobenzofuran]-1,1',3,3'-tetrone, polymer with
     1,4-bis(3,5-diaminobenzoyl)piperazine, 4,4'-oxybis[benzenamine],
     5.5'-oxybis[1.3-isobenzofurandione] and
     4,4'-[1,3-phenylenebis(oxy)]bis[benzenamine] (9CI) (CA INDEX NAME)
     CM
          1
```

CRN 773889-55-7 CMF C18 H22 N6 O2

CM 2

CRN 2479-46-1 CMF C18 H16 N2 O2

CM 3

CRN 2420-87-3 CMF C16 H6 O6

CM 4

CRN 1823-59-2 CMF C16 H6 O7

CM 5

CRN 101-80-4 CMF C12 H12 N2 O

RN 773889-57-9 HCAPLUS CN [5.5'-Biisobenzofura

[5,5'-Biisobenzofuran]-1,1',3,3'-tetrone, polymer with

1,4-bis(3,5-diaminobenzoy1)piperazine,

4,4'-(9H-fluoren-9-ylidene)bis[benzenamine],

5,5'-oxybis[1,3-isobenzofurandione] and 3,3'-sulfonylbis[benzenamine]

(9CI) (CA INDEX NAME)

CM 1

CRN 773889-55-7

CMF C18 H22 N6 O2

CM 2

CRN 15499-84-0 CMF C25 H20 N2

CM 3

CRN 2420-87-3

CMF C16 H6 O6

CM 4

CRN 1823-59-2 CMF C16 H6 O7

CM 5

CRN 599-61-1 CMF C12 H12 N2 O2 S

$$\mathsf{H}_2 \mathsf{N} = \bigcup_{\mathsf{N}} \mathsf{N} \mathsf{H}_2$$

RN 773889-58-0 HCAPLUS

No. (5.5'-Bisobenzofuran]-1,1',3,3'-tetrone, polymer with 3-(4-aminophenoxy)benzenamine, α-[(3-aminopropyl)dimethylsilyl]-ω-[(3-aminopropyl)dimethylsilyl]oxy(dimethylsilylene)] and 1,4-bis(3,5-diaminobenzoyl)piperazine (9CI) (CA INDEX NAME)

CM :

CRN 773889-55-7 CMF C18 H22 N6 O2

CM 2

CRN 97917-34-5

CMF (C2 H6 O Si)n C10 H28 N2 O Si2

CCI PMS

$$H_2N - (CH_2)_3 - si - O - si - O - si - (CH_2)_3 - NH_2$$

CM .

CRN 2657-87-6 CMF C12 H12 N2 O

CM 4

CRN 2420-87-3 CMF C16 H6 O6

RN 773889-62-6 HCAPLUS

CN Benzamide, N,N'-(oxydi-4,1-phenylene)bis[3,5-diamino-, polymer with 3-(4-aminophenoxy)benzenamine, [5,5'-bisobenzofuran]-1,1',3,3'-tetrone, 5,5'-oxybis[1,3-isobenzofurandione] and 4,4'-[1,3-phenylenebis(oxy)]bis[benzenamine] (9CI) (CA INDEX NAME)

CM 1

CRN 773889-59-1 CMF C26 H24 N6 O3

CM 2

CRN 2657-87-6

CMF C12 H12 N2 O

CM 3

CRN 2479-46-1 CMF C18 H16 N2 O2

CM 4

CRN 2420-87-3 CMF C16 H6 O6

CM 5

CRN 1823-59-2 CMF C16 H6 O7

RN 773889-63-7 HCAPLUS

[5,5'-Biisobenzofuran]-1,1',3,3'-tetrone, polymer with 3-(4-aminophenoxy)benzenamine, 5,5'-oxybis[1,3-isobenzofurandione], 4,4'-[1,3-phenylenebis(oxy)]bis[benzenamine] and

5,5'-(sulfonyldi-5,2-benzoxazolediyl)bis[1,3-benzenediamine] (9CI) (CA

19

INDEX NAME)

CM 1

CRN 518992-19-3

CMF C26 H20 N6 O4 S

CM 2

CRN 2657-87-6 CMF C12 H12 N2 O

CM 3

CRN 2479-46-1 CMF C18 H16 N2 O2

CM 4

CRN 2420-87-3

CMF C16 H6 O6

CM 5

CRN 1823-59-2

CMF C16 H6 O7

RN 773889-64-8 HCAPLUS

CN [5,5'-Biisobenzofuran]-1,1',3,3'-tetrone, polymer with

3-(4-aminophenoxy)benzenamine, α -[(3-aminopropyl)dimethylsilyl]-

 $\omega\text{-[[(3-aminopropyl)dimethylsilyl]oxy]poly[oxy(dimethylsilylene)],}\\$

5,5'-oxybis[1,3-isobenzofurandione] and

5,5'-(sulfonyldi-5,2-benzoxazolediyl)bis[1,3-benzenediamine] (9CI) (CA INDEX NAME)

CM 1

CRN 518992-19-3

CMF C26 H20 N6 O4 S

CM 2

CRN 97917-34-5

CMF (C2 H6 O Si)n C10 H28 N2 O Si2

CCI PMS

CM 3

CRN 2657-87-6

CMF C12 H12 N2 O

CM 4

CRN 2420-87-3 CMF C16 H6 O6

CM 5

CRN 1823-59-2 CMF C16 H6 O7

RN 773889-65-9 HCAPLUS

[5,5'-Biisobenzofuran]-1,1',3,3'-tetrone, polymer with 3-(4-aminophenoxy) benzenamine, 9,9-bis(4-aminophenoxy) benzenamine, 2,7-diamine, 5,5'-oxybis[1,3-isobenzofurandione] and 4,4'-[1,3-phenylenebis(oxy)]bis[benzenamine] (9CI) (CA INDEX NAME)

CM :

CN

CRN 176258-99-4 CMF C25 H22 N4

CM 2

CRN 2657-87-6 CMF C12 H12 N2 O

CM 3

CRN 2479-46-1 CMF C18 H16 N2 O2

CM 4

CRN 2420-87-3 CMF C16 H6 O6

CM 5

CRN 1823-59-2 CMF C16 H6 O7

RN 773889-66-0 HCAPLUS

CN 1H,3H-Benzo(1,2-c:4,5-c')difuran-1,3,5,7-tetrone, polymer with [5,5'-blisobenzofuran]-1,1',3,3'-tetrone, 1,4-bis(3,5-diaminobenzoy1)piperazine, 4-methyl-1,3-benzenediamine and

1,4-bis(3,5-diaminobenzoyl)piperazine, 4-methyl-1,3-benzenediamine and
3,3'-[sulfonylbis(4,1-phenyleneoxy)]bis[benzenamine] (9CI) (CA INDEX NAME)

CM 1

CRN 773889-55-7 CMF C18 H22 N6 O2

CM 2

CRN 30203-11-3 CMF C24 H20 N2 O4 S

CM 3

CRN 2420-87-3 CMF C16 H6 O6

CM 4

CRN 95-80-7 CMF C7 H10 N2

CM 5

CRN 89-32-7 CMF C10 H2 O6

RN 773889-67-1 HCAPLUS

[5,5'-Biisobenzofuran]-1,1',3,3'-tetrone, polymer with 1,4-bis(3,5-diaminobenzoyl)piperazine, 5-methyl-1,3-benzenediamine, 5,5'-oxybis[1,3-isobenzofurandione] and 4,4'-[1,3-phenylenebis(oxy)]bis[benzenamine] (9CI) (CA INDEX NAME)

CM 1

CN

CRN 773889-55-7 CMF C18 H22 N6 O2

CM 2

CRN 2479-46-1 CMF C18 H16 N2 O2

CM 3

CRN 2420-87-3 CMF C16 H6 O6

CM 4

CRN 1823-59-2 CMF C16 H6 O7

CM 5

CRN 108-71-4 CMF C7 H10 N2

RN 773889-68-2 HCAPLUS

CN [5,5'-Biisobenzofuran]-1,1',3,3'-tetrone, polymer with 1,4-bis(3,5-diaminobenzoyl)piperazine,

4,4'-(9H-fluoren-9-ylidene)bis[benzenamine], 5,5'-oxybis[1,3-isobenzofurandione] and 4,4'-sulfonylbis[2-aminophenol] (9CI) (CA INDEX NAME)

CM 1

CRN 773889-55-7 CMF C18 H22 N6 O2

CM 2

CRN 15499-84-0 CMF C25 H20 N2

CM 3

CRN 7545-50-8 CMF C12 H12 N2 O4 S

CM ·

CRN 2420-87-3

CMF C16 H6 O6

CM 5

CRN 1823-59-2 CMF C16 H6 O7

RN 773889-69-3 HCAPLUS

CN [5,5'-Biisobenzofuran]-1,1',3,3'-tetrone, polymer with

1,4-bis(3,5-diaminobenzoyl)piperazine,

4,4'-(9H-fluoren-9-ylidene)bis[benzenamine], ar-methyl-1,3-benzenediamine,

5,5'-oxybis[1,3-isobenzofurandione] and

4,4'-[1,3-phenylenebis(oxy)]bis[benzenamine] (9CI) (CA INDEX NAME)

CM 1

CRN 773889-55-7

CMF C18 H22 N6 O2

CM

CRN 26764-44-3

CMF C7 H10 N2

CCI IDS

D1-Me

CM 3

CRN 15499-84-0 CMF C25 H20 N2

CM 4

CRN 2479-46-1 CMF C18 H16 N2 O2

CM 5

CRN 2420-87-3 CMF C16 H6 O6

CM 6

CRN 1823-59-2 CMF C16 H6 O7

RN 773889-70-6 HCAPLUS

CN [5,5'-Biisobenzofuran]-1,1',3,3'-tetrone, polymer with 1,4-bis(3,5-diaminobenzoyl)piperazine, ar-methyl-1,3-benzenediamine, 5,5'-oxybis[1,3-isobenzofurandione] and 4,4'-[1,3-phenylenebis(oxy)]bis[benzenamine] (9CI) (CA INDEX NAME)

CM 1

CRN 773889-55-7 CMF C18 H22 N6 O2

CM 2

CRN 26764-44-3 CMF C7 H10 N2

CCI IDS

D1-Me

CM 3

CRN 2479-46-1 CMF C18 H16 N2 O2

CM 4

CRN 2420-87-3 CMF C16 H6 O6

CM 5

CRN 1823-59-2 CMF C16 H6 O7

RN 773889-71-7 HCAPLUS

CN [5,5'-Biisobenzofuran]-1,1',3,3'-tetrone, polymer with 3-(4-aminophenoxy)benzenamine, 1,4-bis(3,5-diaminobenzoyl)piperazine, 5,5'-oxybis[1,3-isobenzofurandione] and

4,4'-[1,3-phenylenebis(oxy)]bis[benzenamine] (9CI) (CA INDEX NAME)

CM 1

CRN 773889-55-7

CMF C18 H22 N6 O2

CM 2

CRN 2657-87-6 CMF C12 H12 N2 O

CM 3

CRN 2479-46-1 CMF C18 H16 N2 O2

CM 4

CRN 2420-87-3 CMF C16 H6 O6

CM 5

CRN 1823-59-2 CMF C16 H6 O7

RN 773889-72-8 HCAPLUS

CN [5,5'-Biisobenzofuran]-1,1',3,3'-tetrone, polymer with

3-(4-aminophenoxy)benzenamine, α -[(3-aminopropyl)dimethylsilyl]- ω -[[(3-aminopropy1)dimethylsily1]oxy]poly[oxy(dimethylsily1ene)], 1,4-bis(3,5-diaminobenzoy1)piperazine and

5,5'-oxybis[1,3-isobenzofurandione] (9CI) (CA INDEX NAME)

CM

CRN 773889-55-7 CMF C18 H22 N6 O2

CM

CRN 97917-34-5 CMF (C2 H6 O Si)n C10 H28 N2 O Si2 CCI PMS

CM 3

CRN 2657-87-6 CMF C12 H12 N2 O

CM

CRN 2420-87-3 CMF C16 H6 O6

CM 5

CRN 1823-59-2 CMF C16 H6 O7

RN 773889-73-9 HCAPLUS

1H,3H-Benzo[1,2-c:4,5-c']difuran-1,3,5,7-tetrone, polymer with 3-(4-aminophenoxy)benzenamine, [5,5'-biisobenzofuran]-1,1',3,3'-tetrone, 1,4-bis(3,5-diaminobenzoyl)piperazine, 5,5'-oxybis[1,3-isobenzofurandione] and 4,4'-[1,3-phenylenebis(oxy)]bis[benzenamine] (901) (CA INDEX NAME)

CM

CN

CRN 773889-55-7 CMF C18 H22 N6 O2

CM 2

CRN 2657-87-6 CMF C12 H12 N2 O

$$\bigcap_{NH2} \bigcap_{NH2}$$

CM 3

CRN 2479-46-1 CMF C18 H16 N2 O2

CM ·

CRN 2420-87-3 CMF C16 H6 O6

CM 5

CRN 1823-59-2 CMF C16 H6 O7

CM 6

CRN 89-32-7 CMF C10 H2 O6

RN 773889-75-1 HCAPLUS

CN [5,5'-Biisobenzofuran]-1,1',3,3'-tetrone, polymer with

1,4-bis(3,5-diaminobenzoy1)piperazine,

4,4'-(9H-fluoren-9-ylidene)bis[benzenamine], ar-methyl-1,3-benzenediamine,

5,5'-oxybis[1,3-isobenzofurandione] and 4,4'-sulfonylbis[2-aminophenol]

(9CI) (CA INDEX NAME)

CM

CRN 773889-55-7

CMF C18 H22 N6 O2

CM

CRN 26764-44-3

CMF C7 H10 N2

CCI IDS

D1-Me

CM

CRN 15499-84-0 CMF C25 H20 N2

CM

CRN 7545-50-8 CMF C12 H12 N2 O4 S

CM !

CRN 2420-87-3 CMF C16 H6 O6

CM 6

CRN 1823-59-2 CMF C16 H6 O7

RN 773889-76-2 HCAPLUS

CN [5,5'-Biisobenzofuran]-1,1',3,3'-tetrone, polymer with

3-(4-aminophenoxy)benzenamine, α -[(3-aminopropyl)dimethylsilyl]-

ω-[[(3-aminopropyl)dimethylsilyl]oxy]poly[oxy(dimethylsilylene)],

1,4-bis(3,5-diaminobenzoyl)piperazine,

4,4'-(9H-fluoren-9-ylidene)bis[benzenamine], ar-methyl-1,3-benzenediamine and 5,5'-oxybis[1,3-isobenzofurandione] (9CI) (CA INDEX NAME)

CM 1

CRN 773889-55-7

CMF C18 H22 N6 O2

CM 2

CRN 97917-34-5 CMF (C2 H6 O Si)n C10 H28 N2 O Si2 CCI PMS

$$H_{2N-}(CH_{2})_{3-}$$
 S_{1-}^{Me} O_{-} S_{1-}^{Me} O_{-} S_{1-}^{Me} O_{-} $O_{$

CM 3

CRN 26764-44-3 CMF C7 H10 N2 CCI IDS

D1—Me

CM 4

CRN 15499-84-0 CMF C25 H20 N2

CM 5

CRN 2657-87-6 CMF C12 H12 N2 O

CM 6

CRN 2420-87-3 CMF C16 H6 O6

CM 7

CRN 1823-59-2 CMF C16 H6 O7

RN 773889-77-3 HCAPLUS

CN [5,5'-Biisobenzofuran]-1,1',3,3'-tetrone, polymer with 3-(4-aminophenoxy)benzenamine, 1,4-bis(3,5-diaminobenzoy))piperazine, 4,4'-(9H-fluoren-9-ylidene)bis[benzenamine], ar-methyl-1,3-benzenediamine and 5,5'-oxybis[1,3-isobenzofurandione] (9CI) (CA INDEX NAME)

CM

CRN 773889-55-7

CMF C18 H22 N6 O2

CM 2

CRN 26764-44-3 CMF C7 H10 N2 CCI IDS

D1-Me

CM 3

CRN 15499-84-0 CMF C25 H20 N2

CM ·

CRN 2657-87-6 CMF C12 H12 N2 O

$$\bigcap_{NH_2} \bigcap_{NH_2}$$

CM 5

CRN 2420-87-3

CMF C16 H6 O6

CM 6

CRN 1823-59-2 CMF C16 H6 O7

IT 176258-99-4P 518992-19-3P 773889-55-7P

773889-59-1P 773889-61-5P

RL: IMF (Industrial manufacture); RCT (Reactant); PREP (Preparation); RACT (Reactant or reagent)

(monomer; crosslinked polyimides with low dielec. constant, compns. containing them and method for their manufacture and use)

RN 176258-99-4 HCAPLUS

CN 9H-Fluorene-2,7-diamine, 9,9-bis(4-aminophenyl)- (CA INDEX NAME)

RN 518992-19-3 HCAPLUS

CN 1,3-Benzenediamine, 5,5'-(sulfonyldi-5,2-benzoxazolediyl)bis- (9CI) (CA INDEX NAME)

- RN 773889-55-7 HCAPLUS
- CN Methanone, 1,1'-(1,4-piperazinediy1)bis[1-(3,5-diaminopheny1)- (CA INDEX NAME)

- RN 773889-59-1 HCAPLUS
- CN Benzamide, N,N'-(oxydi-4,1-phenylene)bis[3,5-diamino- (9CI) (CA INDEX NAME)

- RN 773889-61-5 HCAPLUS
- CN Benzamide, N, N'-1, 4-phenylenebis [3, 5-diamino- (CA INDEX NAME)

RE.CNT 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT

- L34 ANSWER 4 OF 14 HCAPLUS COPYRIGHT 2009 ACS on STN
- AN 2003:421614 HCAPLUS Full-text
- DN 139:181414
- TI Synthesis and properties of star-like wholly aromatic polyester fibers
- AU Yang, F.; Bai, Y.; Min, B. G.; Kumar, S.; Polk, M. B.
- CS School of Textile and Fiber Engineering, Georgia Institute of Technology, Atlanta, GA, 30332-0295, USA
- SO Polymer (2003), 44(14), 3837-3846 CODEN: POLMAG; ISSN: 0032-3861

- PB Elsevier Science Ltd.
- DT Journal
- LA English
- AB Novel star-like wholly aromatic copolyesters having four arms based on a tetraamine star core, p- and m- hydroxybenzoic acids and 6-hydroxy-2-naphthoic acid have been successfully synthesized and spun into fibers for the investigation of the effect of the star-like structure on improving compressive properties of the fiber. The reactivity of the star core was demonstrated using a model compound with FTIR, 1H and 13C NMR spectroscopy. The 13C NMR spectrum of the star-like terpolymer having a molar ratio of 10:1 of the monomers to star core showed a characteristic peak at around $\delta 62$ ppm which corresponds to a tetra-substituted carbon and thereby demonstrates that the star core was really incorporated into the polymer. The star-like copolyester exhibited a clear stir opalescence and liquid crystalline morphol. in the temperature range of 150-280°. However, no transition was observed in the DSC thermogram except a clear Tq at 110°. The star-like terpolymer fiber, prepared from a polymer with a molar ratio of 500:1 for the monomers to imide core, was spun in the liquid crystalline state at 180°. Fiber structure and properties were studied.
- IT 176258-99-4P
 - RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
- (synthesis and properties of star-like wholly aromatic polyester fibers)
- RN 176258-99-4 HCAPLUS
- CN 9H-Fluorene-2,7-diamine, 9,9-bis(4-aminophenyl)- (CA INDEX NAME)

RE.CNT 24 THERE ARE 24 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT

L34 ANSWER 5 OF 14 HCAPLUS COPYRIGHT 2009 ACS on STN

AN 2001:481874 HCAPLUS Full-text

DN 135:62755

- TI Polyamic acids and polyimides having fluorene structures, and coatings for manufacture of electric insulating films
- IN Matsubara, Minoru; Okada, Takashi; Inoue, Yasutake; Takahashi, Masayuki; Rojanschi, Igor; Goto, Kohei
- PA JSR Co., Ltd., Japan
- SO Jpn. Kokai Tokkyo Koho, 12 pp.

CODEN: JKXXAF

DT Patent

LA Japanese

FAN CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE	
PΙ	JP 2001181390	A	20010703	JP 1999-371369	19991227 <	
PRAI	JP 1999-371369		19991227	<		
OS	MARPAT 135:62755					

OS

AΒ The polyamic acids are reaction products of ≥3-functional polyamines with fluorene structure-containing tetracarboxylic acids I (R1, R2 = H, C1-6 alkyl, C6-14 monocyclic or condensed polycyclic aromatic group) and/or 2,2',3,3'biphenyltetracarboxylic dianhydride. The polyimides are dehydration ringclosure products of the above polyamic acids. The coatings, useful for for interlayer insulating films and elec. insulators in electronic devices, contain the polyamic acids and/or polyimides. Thus, polymerization of 15 mmol 9,9-bis[4-(4-aminophenoxy)phenyl]fluorene and 0.52 mmol 2,7-diamino-9,9-bis[4-(4-aminophenoxy)phenyl]fluorene with 18 mmol 2,2',3,3'-biphenyltetracarboxylic dianhydride at room temperature for 5 h to give a polyamic acid, which was then treated at room temperature for 1 h and then at 100° for 4 h to give a polyimide with dielec. constant 2.95, Tg 326°, solution viscosity (as 10% DMF solution) 53,600 cP, dielec. anisotropy ratio (to the plane and thickness directions) 1.01, and good solubility in N-methylpyrrolidone and cyclohexane. ΙT 346407-43-02

RL: IMF (Industrial manufacture); PRP (Properties); TEM (Technical or engineered material use); PREP (Preparation); USES (Uses)

(polyamic acids and polyimides having fluorene structures for elec. insulating coatings)

RN 346407-43-0 HCAPLUS

CN [4,4'-Biisobenzofuran]-1,1',3,3'-tetrone, polymer with

9,9-bis(4-aminophenyl)-9H-fluorene-2,7-diamine and
4,4'-[9H-fluoren-9-vlidenebis(4,1-phenyleneoxy)]bis[benzenamine] (9CI)

(CA INDEX NAME)

CM

CRN 176258-99-4

CMF C25 H22 N4

CM 2

CRN 47823-88-1

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PAGE 2-A O NH2

CM 3

CRN 3711-04-4 CMF C16 H6 O6

CMF C37 H28 N2 O2

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L34 ANSWER 6 OF 14 HCAPLUS COPYRIGHT 2009 ACS on STN
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AN 2001:143713 HCAPLUS Full-text

DN 134:179710

ΤI Aromatic polymer compositions containing triazene crosslinking agents, their cured films, and manufacture of the films

Akiike, Toshiyuki; Takahashi, Masayuki; Goto, Kohei

IN PA JSR Co., Ltd., Japan

SO Jpn. Kokai Tokkyo Koho, 14 pp.

CODEN: JKXXAF

DT Patent

LA Japanese

FAN.CNT 1

10 / 550887 45 KIND DATE APPLICATION NO. DATE

PI JP 2001055513	A 2001022	7 JP 1999-230444	19990817 <			
PRAI JP 1999-230444	1999081	7 <				
		lator films or optical				
polymers containing	g aromatic ring:	s in structural repeati	ng units, triazenes			
having ≥3 N:NNRR' (R, R' = H, C1-20 alkyl, aryl), and organic solvents. Thus,						
a composition comprising 9,9-bis(4-methylsulfonyloxyphenyl)fluorene-bis(4-						
methylsulfonyloxyphenyl) ether copolymer 100, 2,7-(3,3-dimethyltriazenyl)-9,9-						
bis[4-(3,3- dimethyltriazenyl)phenyl]fluorene 15, and cyclohexanone 900 parts						
		ed at 400° to give a fi				
constant 2.8, elast	tic modulus 5.2	GPa, and good heat and	solvent resistance.			

IT 176258-99-4, 2,7-Diamino-9,9-bis(4-aminophenyl)fluorene RL: RCT (Reactant); RACT (Reactant or reagent)

(in preparation of triazene crosslinking agents for aromatic elec. insulator

optical films)

or

PATENT NO.

RN 176258-99-4 HCAPLUS

CN 9H-Fluorene-2,7-diamine, 9,9-bis(4-aminophenyl)- (CA INDEX NAME)

L34 ANSWER 7 OF 14 HCAPLUS COPYRIGHT 2009 ACS on STN

AN 2001:143649 HCAPLUS Full-text

DN 134:179368

TI Tetrakistriazenyl fluorene compounds for polymer crosslinking agents

IN Akiike, Toshiyuki; Goto, Kohei

PA JSR Co., Ltd., Japan

SO Jpn. Kokai Tokkyo Koho, 9 pp.

CODEN: JKXXAF

DT Patent

LA Japanese

FAN.CNT 1

PATENT NO. KIND DATE APPLICATION NO. DATE ----JP 1999-230443 PI JP 2001055368 A 20010227 19990817 <--PRAI JP 1999-230443 19990817 <--

GI

OS MARPAT 134:179368

- AB The triazenyl-containing fluorene compds. I (R1, R2 = H, C1-20-alkyl, aryl, at Relation and the compds of the c
- IT 176258-99-4, 2,7-Diamino-9,9-bis(4-aminophenyl)fluorene RL: RCT (Reactant); RACT (Reactant or reagent)

(tetrakistriazenyl fluorene compds. for polymer crosslinking agents)
RN 176258-99-4 HCAPLUS

CN 9H-Fluorene-2,7-diamine, 9,9-bis(4-aminophenyl)- (CA INDEX NAME)

- L34 ANSWER 8 OF 14 HCAPLUS COPYRIGHT 2009 ACS on STN
- AN 1999:788218 HCAPLUS Full-text
- DN 132:37035
- TI Thermosetting resin compositions and their curing materials having

excellent storage stability, transparency, adhesion, electric insulating property, and crack, heat, and moisture resistance

- IN Matsubara, Minoru; Inoue, Yasutake; Kakuta, Mayumi; Goto, Kohei; Kurosawa, Takahiko; Shinoda, Tomotaka; Yamada, Kinji
- PA JSR Co., Ltd., Japan
- SO Jpn. Kokai Tokkyo Koho, 19 pp.
- CODEN: JKXXAF
- DT Patent LA Japanese
- TH Oabane

FAN.CNT 1					
PATENT NO.	KIND	DATE	APPLICATION NO.	DATE	
PI JP 11343414	A	19991214	JP 1998-164447	19980529 <	
PRAI JP 1998-164447		19980529	<		

AB Title compns., useful for coatings, comprise hydrolyzed organosilanes and/or partially condensed hydrolyzable organosilanes and polyimides having ≥ 1

10 / 550887

reactive group and/or their polyamic acids. Polyamides are manufactured from compds. containing ≥ 3 amino groups, carboxylic acid dianhydrides, and diamines. Thus, a coating containing a polyimide (prepared from $9,9-\mathrm{bis}\,[4-(4-\mathrm{aminophenoxy})\mathrm{phenyl}]\mathrm{fluorene}$ 5,327, 2,2-bis(3,4-dicarboxyphenyl)-1,1,1,3,3,3-hexafluoropropane dianhydride 6.078, and 2,7-diamino-9,9-bis[4-(4-\mathrm{aminophenoxy})\mathrm{phenyl}]\mathrm{fluorene} 0.422 g) 1.57, di-

isopropoxybis(ethylacetoacetate)titanium 2.63, and MeSiOMe3 6.36 g showed 5% weight-loss temperature $>600^{\circ}$ and dielec. constant 2.3.

252370-79-9P 252370-80-2P 252370-81-3P

RL: IMF (Industrial manufacture); PRP (Properties); TEM (Technical or engineered material use); PREP (Preparation); USES (Uses)

(thermosetting resin compns. having good storage stability, transparency, adhesion, elec. insulating property, and crack, heat, and

moisture resistance)

RN 252370-79-9 HCAPLUS CN 1,3-Isobenzofurandione, 5,5'-[2,2,2-trifluoro-1-

(trifluoromethyl)ethylidenelbis-, polymer with

9,9-bis(4-aminophenyl)-9H-fluorene-2,7-diamine and

 ${\tt 4,4'-[9H-fluoren-9-ylidenebis(4,1-phenyleneoxy)]bis[benzenamine]} \eqno(9CI)$

(CA INDEX NAME)

CM

ТТ

CRN 176258-99-4 CMF C25 H22 N4

CM 2

CRN 47823-88-1

CMF C37 H28 N2 O2

CM 3

CRN 1107-00-2 CMF C19 H6 F6 O6

RN 252370-80-2 HCAPLUS

[4,4'-Biisobenzofuran]-1,1',3,3'-tetrone, polymer with 9,9-bis(4-aminophenyl)-9H-fluorene-2,7-diamine and 4,4'-(9H-fluoren-9-ylidene)bis(benzenamine) (9CI) (CA INDEX NAME)

CM 1

CN

CRN 176258-99-4 CMF C25 H22 N4

CM 2

CRN 15499-84-0 CMF C25 H20 N2

CM

CRN 3711-04-4 CMF C16 H6 O6

RN 252370-81-3 HCAPLUS

CN Benzoic acid, 3,5-diamino-, polymer with [4,4'-biisobenzofuran]-1,1',3,3'-tetrone, 9,9-bis(4-aminophenyl)-9H-fluorene-2,7-diamine and 4,4'-(9H-fluoren-9-ylidene)bis(benzenamine) (9CI) (CA INDEX NAME)

CM 1

CRN 176258-99-4 CMF C25 H22 N4

CM :

CRN 15499-84-0

10 / 550887 50

CMF C25 H20 N2

CM

CRN 3711-04-4 CMF C16 H6 O6

CM

CRN 535-87-5 CMF C7 H8 N2 O2

- L34 ANSWER 9 OF 14 HCAPLUS COPYRIGHT 2009 ACS on STN
- 1999:310082 HCAPLUS Full-text AN
- DN 131:102606
- Synthesis and characterization of high performance polymers containing fluorene central units
- AU Harruna, Issifu I.; Petzold, Odessa N.; Bray, Melody L.
- CS Department of Chemistry and the High Performance Polymers and Composites Center, Clark Atlanta University, Atlanta, GA, 30314, USA

- SO Recent Research Developments in Polymer Science (1998), 2(Pt. 2), 183-201 CODEN. RNDPFX
- PB Transworld Research Network
- DT Journal

AB

- LA English
 - Radially oriented polyimide, polybenzoxazole (PBO), and polybenzothiazole (PBT) systems were prepared based on the geometrically favorable fluorene central unit. The three-dimensional central unit, 2,7-diamino-9,9-bis(4aminophenvl)fluorene, was used to effectively control the orientation of linear polymer chains. The synthesis of the central fluorene unit involved the reaction of 2.7-dinitro-9-fluorenone with aniline in the presence of aniline hydrochloride followed by the reduction of the nitro functionalities. Polyimide systems exhibiting improved solubility in organic solvents and strong acids, transitions at lower temps. and good thermal properties were prepared by introducing hexafluoro groups (bis-4,4'aminophenylhexafluoropropane and 4,4'-(hexafluoroisopropylidene)diphthalic anhydride), and by using a dianhydride (bicyclo[2.2.2]oct-7-ene-2,3,5,6tetracarboxylic dianhydride) with reduced symmetry and decreased π -electron d. relative to benzene tetracarboxylic dianhydride. An improvement in the solubility of the thermally stable radially oriented polyimides was observed as a result of a decrease in linearity due to the radial structure of the three-dimensional polyimides. Polyimides containing the hexafluoroisopropyl group and the bicyclic ring (bicyclo[2.2.2]oct-7-ene-2,3,5,6-tetracarboxylic dianhydride) were prepared to increase solubility while maintaining thermal stability. The hexafluoroisopropylidene group was used to increase chain flexibility while the presence of fluoro groups led to improved solubility without compromising thermal stability. Bicvclo[2.2.2]oct-7-ene-2,3,5,6tetracarboxylic dianhydride was used to enhance the processability of the polyimides by improving tractability. The radially oriented polymers are expected to be excellent candidates of improved compressive properties compared to the linear analogs. The polybenzoxazoles (PBOs and PBTs) are highly solvent resistant, and exhibit high thermal and thermooxidative stabilities. The C9 of the central units of the polymers showed 13C NMR
- RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(core; preparation of chemical resistant thermally stable radial polyimides and

- polybenzoxazoles and polybenzothiazoles with aminophenyl-fluorene central units)
- RN 176258-99-4 HCAPLUS
- CN 9H-Fluorene-2,7-diamine, 9,9-bis(4-aminophenyl)- (CA INDEX NAME)

chemical shifts in the δ 64.7-73.7 ppm region. 176258-99-49, 2.7-Diamino-9.9-bis(4-aminophenyl)fluorene

RE.CNT 15 THERE ARE 15 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

52

- L34 ANSWER 10 OF 14 HCAPLUS COPYRIGHT 2009 ACS on STN
- AN 1999:59121 HCAPLUS Full-text
- DN 130:125670
- TI Thermal and nuclear magnetic resonance characterization of linear and radial high-performance polyimides
- AU Harruna, Issifu I.; Petzold, Odessa
- CS Dep. Chem., High-Performance Polymers Composites Center, Clark Univ., Atlanta, GA, 30314, USA
- SO International Journal of Polymer Analysis and Characterization (1998), 4(6), 565-578
 - CODEN: IPACEZ; ISSN: 1023-666X
- PB Gordon & Breach Science Publishers
- DT Journal
- LA English
- Radially oriented polyimides were prepared based on the geometrically favorable fluorene central unit by condensation with the C-terminus of the linear polyimides. Characterization of the polymers was achieved by Fourier-transform IR resonance (FTIR) spectroscopy, 1H NMR and 13C NMR spectroscopy, x-ray diffraction, solubility, and inherent viscosity. Thermal properties were investigated by thermogravimetric anal. (TGA), differential calorimetry scanning (DSC), and optical polarizing microscopy. The radially oriented polyimides were compared to their linear analogs. Thermogravimetric analyses indicate that linear and radially oriented polyimides were stable .ltorsim.400° in air and under Ar atmospheres.
- RN 176258-99-4 HCAPLUS
- CN 9H-Fluorene-2.7-diamine, 9.9-bis(4-aminophenyl)- (CA INDEX NAME)

- L34 ANSWER 11 OF 14 HCAPLUS COPYRIGHT 2009 ACS on STN
- AN 1997:388732 HCAPLUS Full-text
- DN 127:26374
- OREF 127:4979a,4982a
- TI Polyamic acid or polyimide liquid crystal-orienting agent
- IN Nishikawa, Michinori; Kawamura, Shigeo; Toyoshima, Hitoshi; Matsuki, Yasuo
- PA Japan Synthetic Rubber Co., Ltd., Japan
- SO Jpn. Kokai Tokkyo Koho, 15 pp.
- CODEN: JKXXAF
- DT Patent
- LA Japanese
- FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	JP 09090367	A	19970404	JP 1995-243041	19950921 <
	JP 3203634	B2	20010827		
PRAI	JP 1995-243041		19950921	<	

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

AB The liquid crystal-orienting agent is a polyamic acid obtained by treating a tetracarboxylic dianhydride I with (A) a phenylenediamine II, (B) ≥1 aromatic diamine selected from III-VII, and (C) a steroid structure-having diamine H2NR14NH2 (R1 = tetravalent organic group; R2-6, R8-13 = alkyl, alkoxy, halo; a = 0-4; R7 = divalent organic group; b, c, h, i, j = 0-3; d, e, f, g, k = 0-4; R14 = steroid structure-having divalent organic group) or a polyamide obtained by dehydration cyclization of thus obtained polyamic acid. Liquidcrystal oriented films containing the orienting agent show uniform thickness, good crystal orientation, and long reliability.

ΙT 190196-08-89

RL: DEV (Device component use); IMF (Industrial manufacture); TEM (Technical or engineered material use); PREP (Preparation); USES (Uses) (polyamic acid or polyimide liquid crystal-orienting agent showing uniform thickness)

190196-08-8 HCAPLUS RN

Cholest-5-en-3-ol (3β) -, 3,5-diaminobenzoate, polymer with CN 1,4-benzenediamine, 9,9-bis(4-aminophenyl)-9H-fluorene-2,7-diamine and hexahydrofuro[3',4':4,5]cyclopenta[1,2-c]pyran-1,3,4,6-tetrone (9CI) (CA INDEX NAME)

CM 1

CRN 176258-99-4 CMF C25 H22 N4

CM 2

CRN 173027-19-5 CMF C34 H52 N2 O2

Absolute stereochemistry.

10 / 550887 54

CM

CRN 127804-19-7 CMF C10 H8 O6

CM

CRN 106-50-3 CMF C6 H8 N2

L34 ANSWER 12 OF 14 HCAPLUS COPYRIGHT 2009 ACS on STN

AN 1997:252345 HCAPLUS Full-text

126:277826

OREF 126:53875a,53878a

ΤI Synthesis of three-dimensional polyimides based on the fluorene unit for compressive strength studies

AII Petzold, Odessa N.; Harruna, Issifu J.; Bota, Kofi B.; Dean, Derrick R. CS Department of Chemistry and the High Performance Polymers and Composites

Center, Clark Atlanta University, Atlanta, GA, 30314, USA

SO Polymeric Materials Science and Engineering (1997), 76, 179-180

CODEN: PMSEDG; ISSN: 0743-0515

PB American Chemical Society English

DT Journal

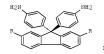
LA

- AB The preparation of star-like polyimides by by condensing linear polyimides with 2,7-diamino-9,9-bis(4-aminophenyl)fluorene is reported. The properties of these 3-dimensional polyimides are discussed.
- IT 176258-99-4DP, polyimides
 - RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation) (preparation and properties of star-like polyimides)
- RN 176258-99-4 HCAPLUS
- CN 9H-Fluorene-2,7-diamine, 9,9-bis(4-aminophenyl)- (CA INDEX NAME)

- L34 ANSWER 13 OF 14 HCAPLUS COPYRIGHT 2009 ACS on STN
- AN 1996:242795 HCAPLUS Full-text
- DN 124:316706

OREF 124:58733a,58736a

- TI Convenient synthesis of 2,7-diamino-9,9-bis(4-aminophenyl)fluorene
- AU Lian, Guohua; Polk, Malcolm B.
- CS School Textile and Fiber Engineering, Georgia Inst. Technol., Atlanta, GA, 30332, USA
- SO Synthetic Communications (1996), 26(10), 2031-6
- CODEN: SYNCAV; ISSN: 0039-7911
- PB Dekker DT Journal
- LA English
- OS CASREACT 124:316706
- GI



- AB For the synthesis of the title compound (I, R = NH2), first I (R = NO2) was synthesized by the reaction of 2,7-dinitro-9-fluorenone with aniline and aniline hydrochloride. I (R = NH2) was obtained by the reduction of I (R = NO2) with hydrazine hydrate and 10% palladium on carbon.
 - I 176258-99-4P

RL: SPN (Synthetic preparation); PREP (Preparation) (synthesis of diaminobis(aminophenyl)fluorene)

- RN 176258-99-4 HCAPLUS
- CN 9H-Fluorene-2,7-diamine, 9,9-bis(4-aminophenyl)- (CA INDEX NAME)

L34 ANSWER 14 OF 14 HCAPLUS COPYRIGHT 2009 ACS on STN

1995:817796 HCAPLUS Full-text AN

DN 123:229141

OREF 123:40953a,40956a

Aryl fluoride monomers in nucleophilic aromatic substitution

polymerization: evaluation of monomer reactivity by 19F NMR spectroscopy

AII Carter, Kenneth R. CS

IBM Res. Div., Almaden Res. Cent., San Jose, CA, 95120-6099, USA

SO Macromolecules (1995), 28(19), 6462-70 CODEN: MAMOBX; ISSN: 0024-9297

PB American Chemical Society

DT Journal

LA English AB

The reactivity of a number of aryl fluoride monomers used in nucleophilic aromatic substitution polymerization was explored utilizing 19F NMR expts. NMR is a valuable tool for evaluating the electron-withdrawing effect of substituents present on Ph rings. When an electron-withdrawing group is present on a Ph ring, a partial pos. charge develops at the ortho and para positions through resonance interactions. While both 13C and 19F NMR were used to probe the electron d. at the actual site of nucleophilic reaction, 19F NMR chemical shifts proved to be the most sensitive probe, with a chemical shift range spanning 2500 Hz between the most activated monomer examined, difluorodiphenyl sulfone, and nonactivated fluorobenzene. The 19F shifts reflect the reactivity of the individual monomers examined Taft inductive and resonance parameters were calculated for a series of monomers from 19F data and used to identify activating forces for the monomers. NMR data were compared with calculated net atomic charges. Relative reactivity studies were also performed in order to verify the utility of this fast and convenient NMR probe of monomer reactivity.

168914-99-6

RL: PRP (Properties); RCT (Reactant); RACT (Reactant or reagent) (evaluation of reactivity of aryl fluoride monomers in nucleophilic aromatic substitution polymerization)

RN 168914-99-6 HCAPLUS

CN 5,6'-Bibenzoxazole, 2,2'-bis(4-fluorophenvl)- (CA INDEX NAME)

=> d his

SET COST OFF

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